

V. Balasubramanian

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LOGINID:ssspal611bxv

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

* * * * * Welcome to STN International * * * * *

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
NEWS 2 "Ask CAS" for self-help around the clock
NEWS 3 May 12 EXTEND option available in structure searching
NEWS 4 May 12 Polymer links for the POLYLINK command completed in REGISTRY
NEWS 5 May 27 New UPM (Update Code Maximum) field for more efficient patent
SDIs in CAPLUS
NEWS 6 May 27 CAPLUS super roles and document types searchable in REGISTRY
NEWS 7 Jun 28 Additional enzyme-catalyzed reactions added to CASREACT
NEWS 8 Jun 28 ANTE, AQUALINE, BIOENG, CIVILENG, ENVIROENG, MECHENG,
and WATER from CSA now available on STN(R)
NEWS 9 Jul 12 BEILSTEIN enhanced with new display and select options,
resulting in a closer connection to BABS
NEWS 10 Jul 30 BEILSTEIN on STN workshop to be held August 24 in conjunction
with the 228th ACS National Meeting
NEWS 11 AUG 02 IFIPAT/IFIUDB/IFICDB reloaded with new search and display
fields
NEWS 12 AUG 02 CAPLUS and CA patent records enhanced with European and Japan
Patent Office Classifications
NEWS 13 AUG 02 STN User Update to be held August 22 in conjunction with the
228th ACS National Meeting
NEWS 14 AUG 02 The Analysis Edition of STN Express with Discover!
(Version 7.01 for Windows) now available
NEWS 15 AUG 04 Pricing for the Save Answers for SciFinder Wizard within
STN Express with Discover! will change September 1, 2004

NEWS EXPRESS JULY 30 CURRENT WINDOWS VERSION IS V7.01, -CURRENT
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
AND CURRENT DISCOVER FILE IS DATED 11 AUGUST 2004
NEWS HOURS STN Operating Hours Plus Help Desk Availability
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* * * * * STN Columbus * * * * *

10/725,079

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FILE 'HOME' ENTERED AT 16:38:28 ON 17 AUG 2004

=> file reg

COST IN U.S. DOLLARS

SINCE FILE

TOTAL

ENTRY

SESSION

FULL ESTIMATED COST

0.21

0.21

FILE 'REGISTRY' ENTERED AT 16:38:38 ON 17 AUG 2004

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 16 AUG 2004 HIGHEST RN 727651-15-2

DICTIONARY FILE UPDATES: 16 AUG 2004 HIGHEST RN 727651-15-2

TSCA INFORMATION NOW CURRENT THROUGH MAY 21, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:

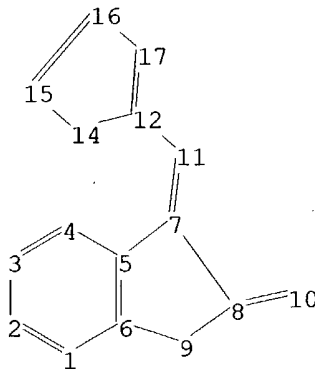
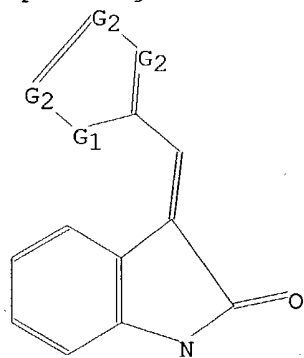
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=>

Uploading C:\STNEXP4\QUERIES\107250792.str



chain nodes :

10 11

ring nodes :

1 2 3 4 5 6 7 8 9 12 14 15 16 17

chain bonds :

7-11 8-10 11-12

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ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 12-14 12-17 14-15 15-16 16-17

exact/norm bonds :

5-7 6-9 7-8 7-11 8-9 8-10 11-12 12-14 12-17 14-15 15-16 16-17

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6

isolated ring systems :

containing 1 : 12 :

G1:O,S,N

G2:C,O,S,N

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:CLASS

11:CLASS 12:Atom 14:Atom 15:Atom 16:Atom 17:Atom

L1 STRUCTURE UPLOADED

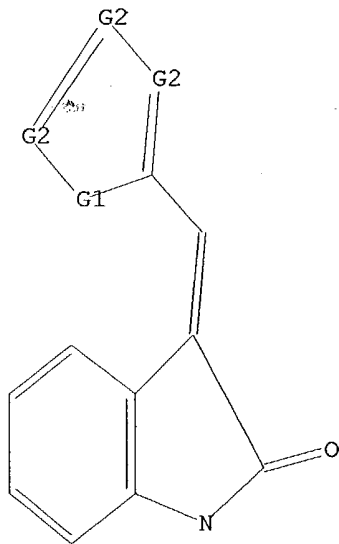
=> que L1

L2 QUE L1

=> d l1

L1 HAS NO ANSWERS

L1 STR



G1 O,S,N

G2 C,O,S,N

Structure attributes must be viewed using STN Express query preparation.

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=> s ll sss sam

SAMPLE SEARCH INITIATED 16:39:00 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 634 TO ITERATE

100.0% PROCESSED 634 ITERATIONS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

50 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 11170 TO 14190
PROJECTED ANSWERS: 2371 TO 3869

L3 50 SEA SSS SAM L1

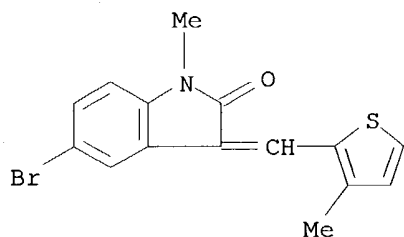
=> d scan

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 2H-Indol-2-one, 5-bromo-1,3-dihydro-1-methyl-3-[(3-methyl-2-thienyl)methylene]- (9CI)

MF C15 H12 Br N O S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

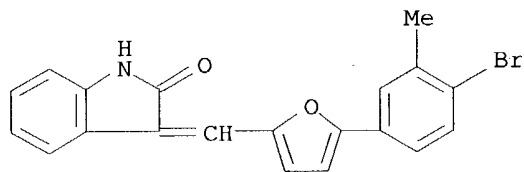
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):49

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 2H-Indol-2-one, 3-[[5-(4-bromo-3-methylphenyl)-2-furanyl]methylene]-1,3-dihydro- (9CI)

MF C20 H14 Br N O2



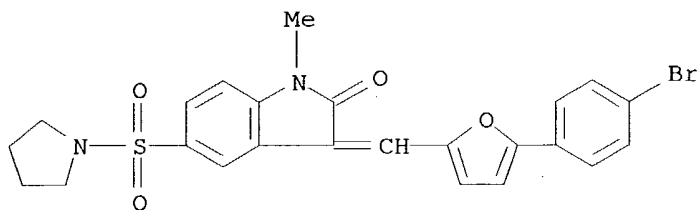
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Pyrrolidine, 1-[[3-[[5-(4-bromophenyl)-2-furanyl]methylene]-2,3-dihydro-1-methyl-2-oxo-1H-indol-5-yl]sulfonyl]- (9CI)

MF C24 H21 Br N2 O4 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

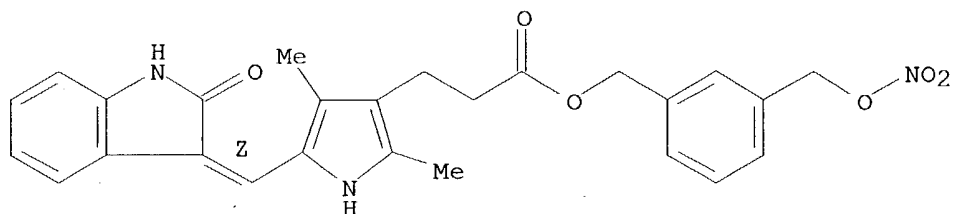
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Pyrrole-3-propanoic acid, 5-[(Z)-(1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2,4-dimethyl-, [3-[(nitrooxy)methyl]phenyl]methyl ester (9CI)

MF C26 H25 N3 O6

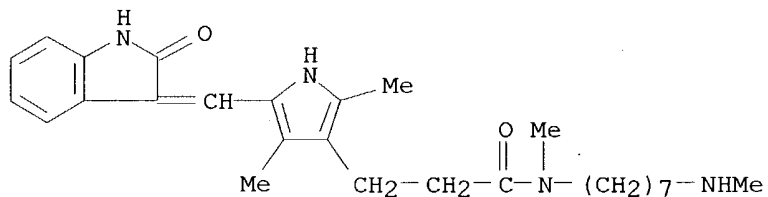
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-propanamide, 5-[(1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-
N,2,4-trimethyl-N-[7-(methylamino)heptyl]- (9CI)
MF C27 H38 N4 O2
CI COM

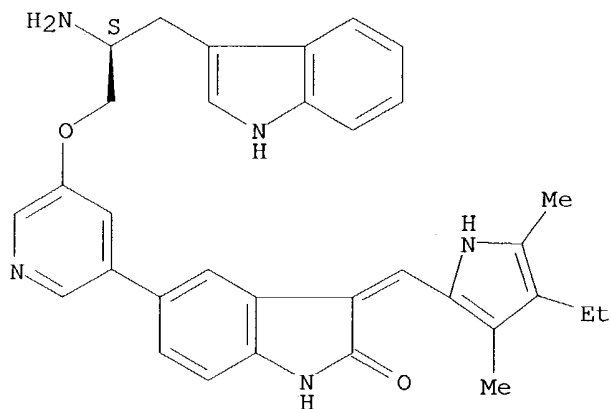


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 5-[5-[(2S)-2-amino-3-(1H-indol-3-yl)propoxy]-3-pyridinyl]-
3-[(4-ethyl-3,5-dimethyl-1H-pyrrol-2-yl)methylene]-1,3-dihydro- (9CI)
MF C33 H33 N5 O2
CI COM

Absolute stereochemistry.
Double bond geometry unknown.

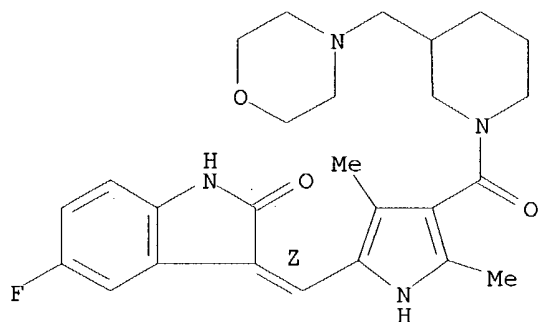


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperidine, 1-[[5-[(Z)-(5-fluoro-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-3-(4-morpholinylmethyl)- (9CI)
MF C26 H31 F N4 O3

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

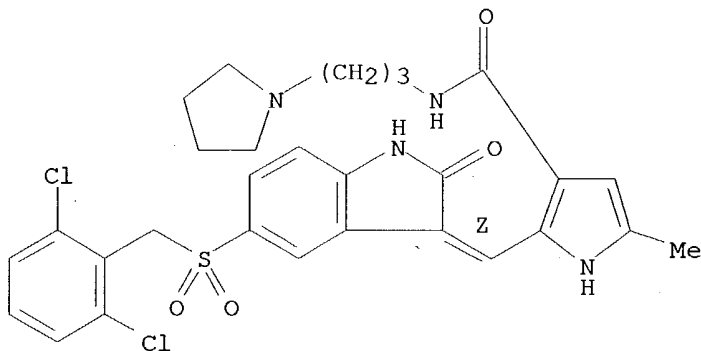
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Pyrrole-3-carboxamide, 2-[(Z)-[5-[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-5-methyl-N-[3-(1-pyrrolidinyl)propyl]- (9CI)

MF C29 H30 Cl2 N4 O4 S

Double bond geometry as shown.

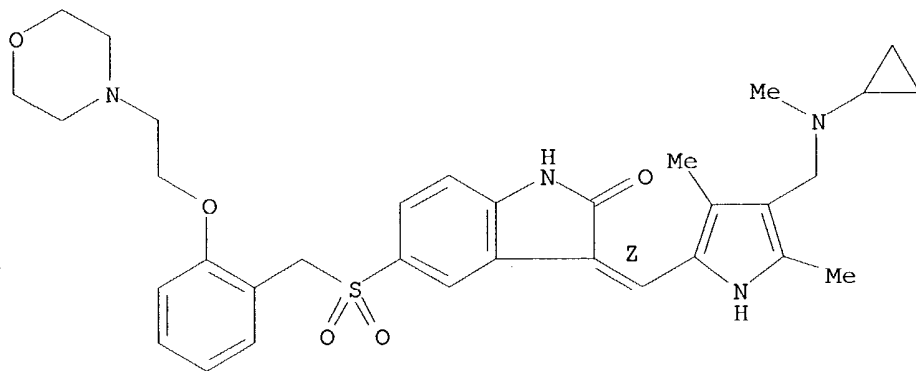


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 3-[[4-[(cyclopropylmethylamino)methyl]-3,5-dimethyl-1H-
pyrrol-2-yl]methylene]-1,3-dihydro-5-[[[2-[2-(4-
morpholinyl)ethoxy]phenyl]methyl]sulfonyl]-, (3Z)- (9CI)
MF C33 H40 N4 O5 S

Double bond geometry as shown.

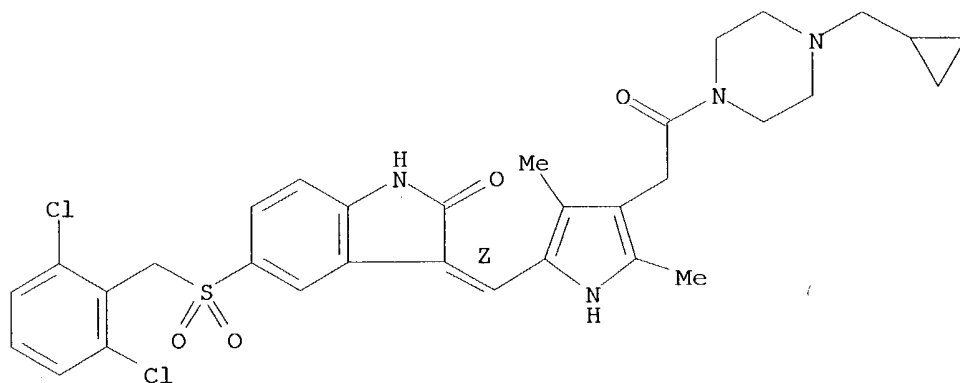


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-(cyclopropylmethyl)-4-[[5-[(Z)-[5-[(2,6-
dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-
ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]acetyl]- (9CI)
MF C32 H34 Cl2 N4 O4 S

Double bond geometry as shown.

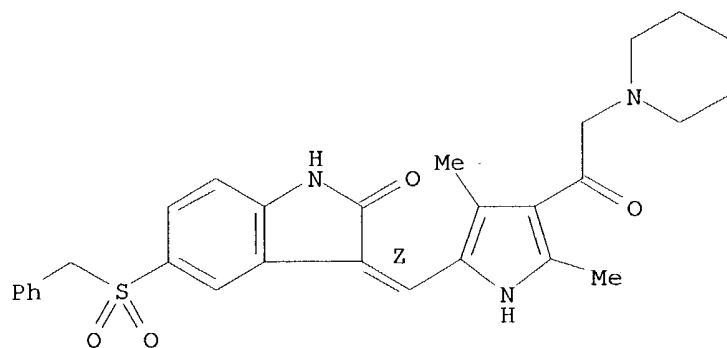


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 3-[[3,5-dimethyl-4-(1-piperidinylacetyl)-1H-pyrrol-2-yl]methylene]-1,3-dihydro-5-[(phenylmethyl)sulfonyl]-, (3Z)- (9CI)
MF C29 H31 N3 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

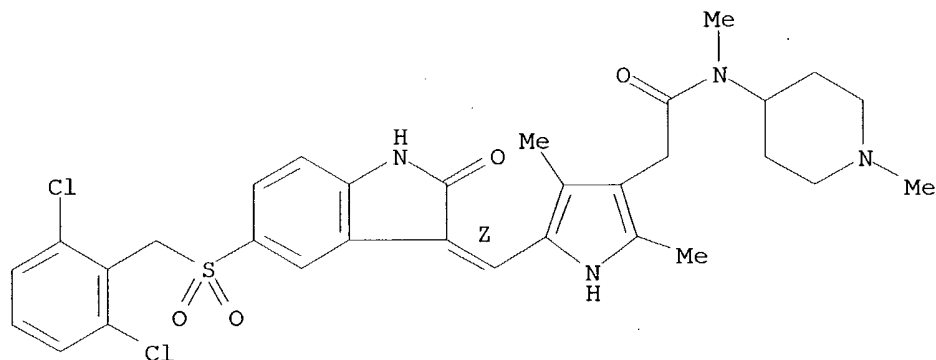
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Pyrrole-3-acetamide, 5-[(Z)-[5-[[[(2,6-dichlorophenyl)methyl]sulfonyl]-
1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-N,2,4-trimethyl-N-(1-methyl-4-
piperidiny)]- (9CI)

MF C31 H34 Cl2 N4 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

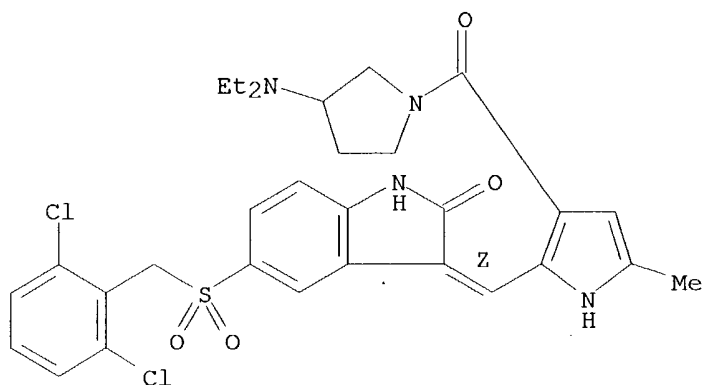
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 3-Pyrrolidinamine, 1-[[2-[(Z)-[5-[[[(2,6-dichlorophenyl)methyl]sulfonyl]-
1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-5-methyl-1H-pyrrol-3-
yl]carbonyl]-N,N-diethyl- (9CI)

MF C30 H32 Cl2 N4 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

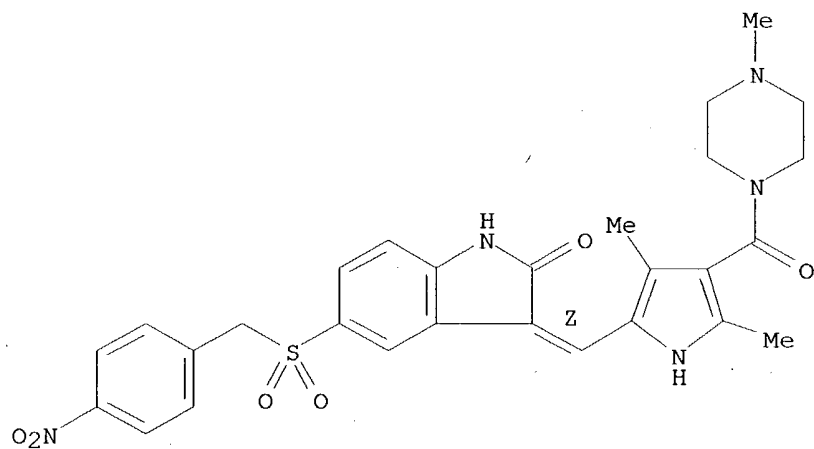
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Piperazine, 1-[[5-[(Z)-[1,2-dihydro-5-[[4-nitrophenyl)methyl]sulfonyl]-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-4-methyl- (9CI)

MF C28 H29 N5 O6 S

Double bond geometry as shown.

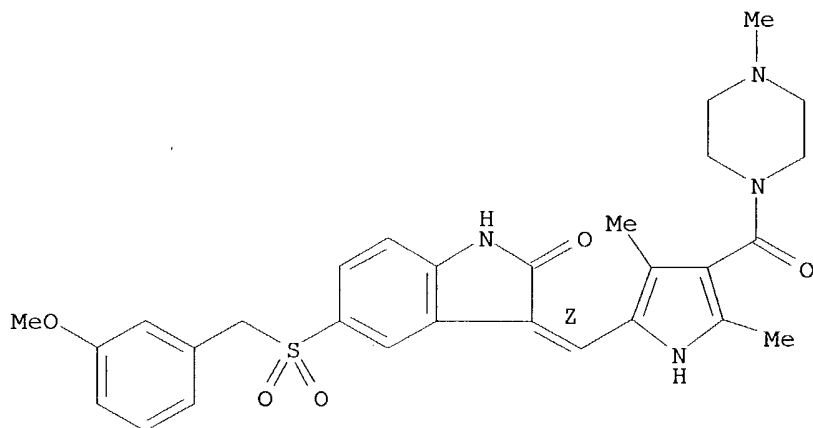


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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[(Z)-[1,2-dihydro-5-[[3-methoxyphenyl)methyl]sulfonyl]-
2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-4-
methyl- (9CI)
MF C29 H32 N4 O5 S

Double bond geometry as shown.



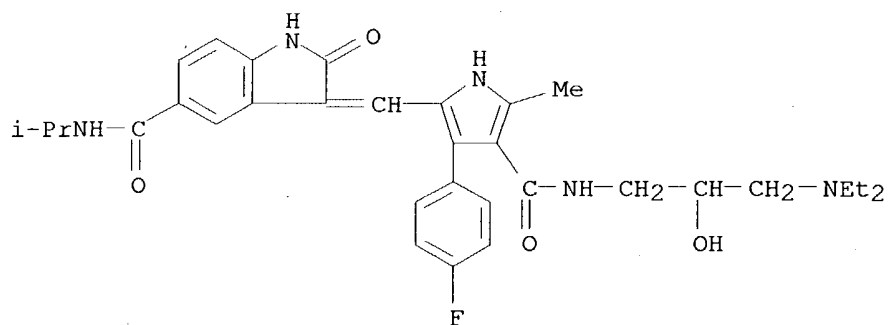
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Indole-5-carboxamide, 3-[[4-[[[3-(diethylamino)-2-hydroxypropyl]amino]carbonyl]-3-(4-fluorophenyl)-5-methyl-1H-pyrrol-2-yl]methylene]-2,3-dihydro-N-(1-methylethyl)-2-oxo- (9CI)

MF C32 H38 F N5 O4

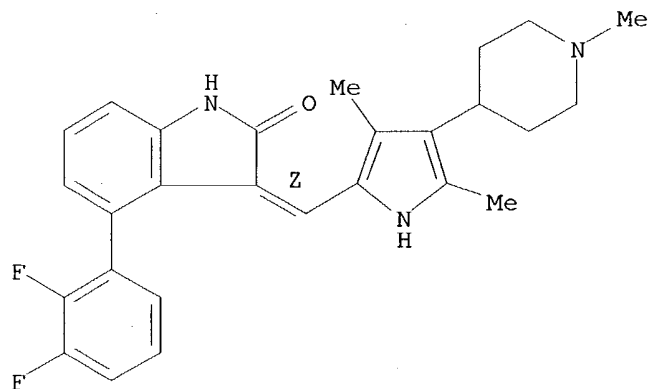


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 4-(2,3-difluorophenyl)-3-[[3,5-dimethyl-4-(1-methyl-4-
piperidinyl)-1H-pyrrol-2-yl]methylene]-1,3-dihydro-, (3Z)- (9CI)
MF C27 H27 F2 N3 O

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

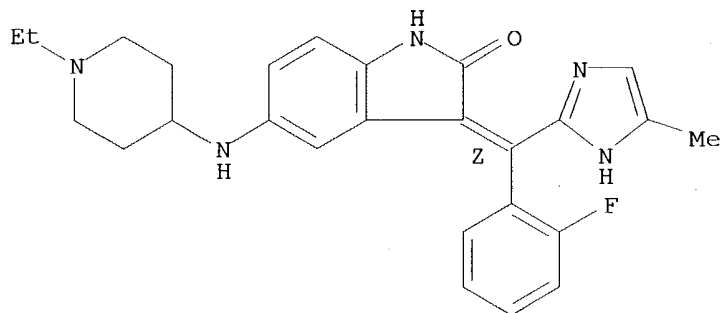
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 2H-Indol-2-one, 5-[(1-ethyl-4-piperidinyl)amino]-3-[(2-fluorophenyl)(4-methyl-1H-imidazol-2-yl)methylene]-1,3-dihydro-, (3Z)- (9CI)

MF C26 H28 F N5 O

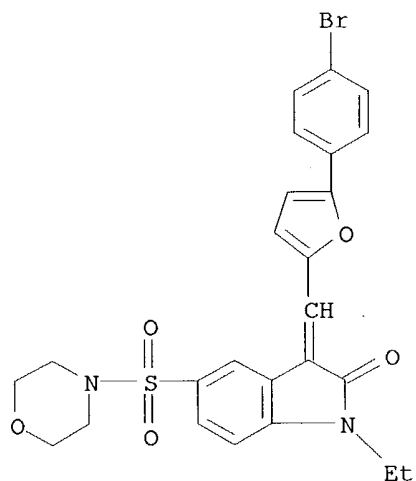
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Morpholine, 4-[[3-[[5-(4-bromophenyl)-2-furanyl]methylene]-1-ethyl-2,3-
dihydro-2-oxo-1H-indol-5-yl]sulfonyl]- (9CI)
MF C25 H23 Br N2 O5 S



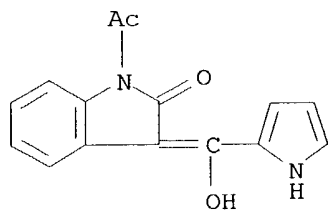
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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IN 2H-Indol-2-one, 1-acetyl-1,3-dihydro-3-(hydroxy-1H-pyrrol-2-ylmethylene)-
(9CI)

MF C15 H12 N2 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

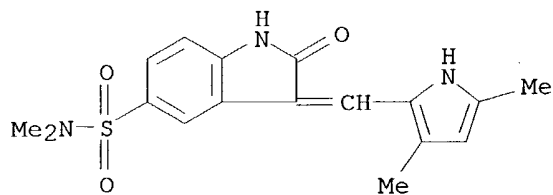
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IN 1H-Indole-5-sulfonamide, 3-[(3,5-dimethyl-1H-pyrrol-2-yl)methylene]-2,3-

dihydro-N,N-dimethyl-2-oxo- (9CI)

MF C17 H19 N3 O3 S



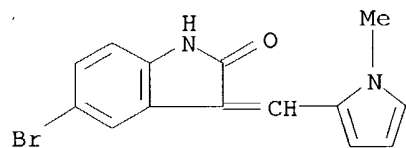
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 2H-Indol-2-one, 5-bromo-1,3-dihydro-3-[(1-methyl-1H-pyrrol-2-yl)methylene]-
(9CI)

MF C14 H11 Br N2 O



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

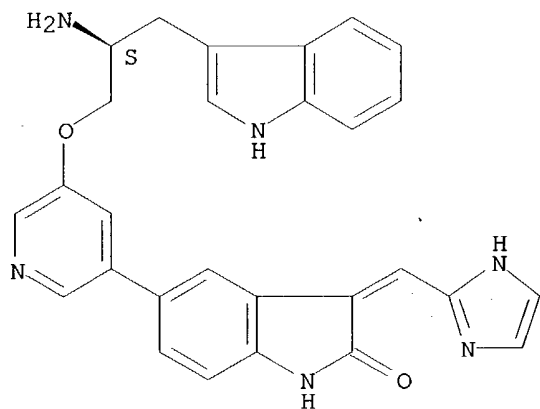
IN 2H-Indol-2-one, 5-[5-[(2S)-2-amino-3-(1H-indol-3-yl)propoxy]-3-pyridinyl]-
1,3-dihydro-3-(1H-imidazol-2-ylmethylene)- (9CI)

MF C28 H24 N6 O2

CI COM

Absolute stereochemistry.

Double bond geometry unknown.

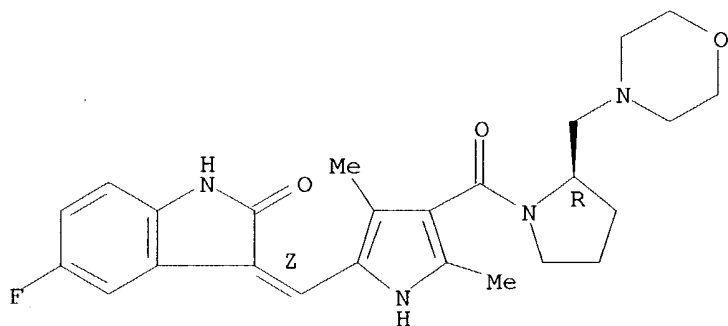


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Pyrrolidine, 1-[[5-[(Z)-(5-fluoro-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-2-(4-morpholinymethyl)-, (2R)- (9CI)
MF C25 H29 F N4 O3

Absolute stereochemistry.
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

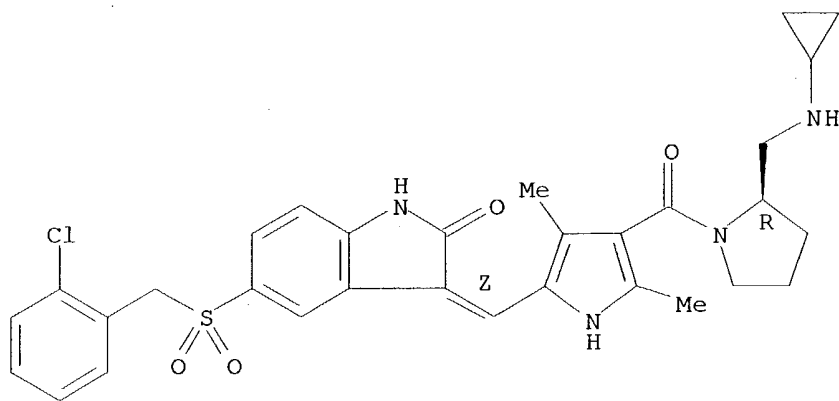
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 2-Pyrrolidinemethanamine, 1-[[5-[(Z)-[5-[(2-chlorophenyl)methyl]sulfonyl]-
1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-
yl]carbonyl]-N-cyclopropyl-, (2R)- (9CI)

MF C31 H33 Cl N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

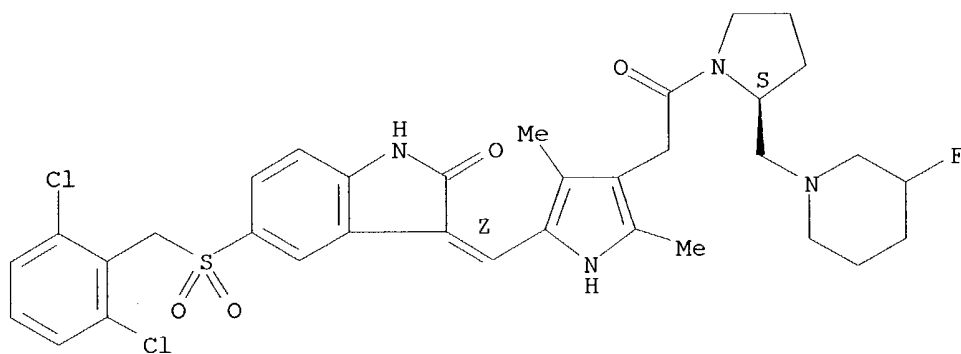
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Pyrrolidine, 1-[[5-[(Z)-[5-[[2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]acetyl]-2-[(3-fluoro-1-piperidinyl)methyl]-, (2S)- (9CI)

MF C34 H37 Cl2 F N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

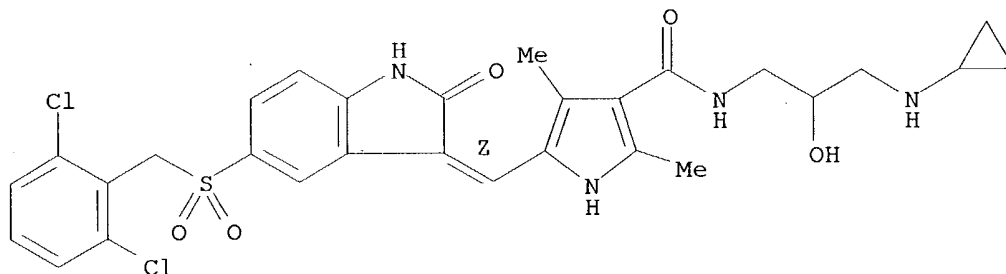
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Pyrrole-3-carboxamide, N-[3-(cyclopropylamino)-2-hydroxypropyl]-5-[(Z)-
[5-[[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-
ylidene]methyl]-2,4-dimethyl- (9CI)

MF C29 H30 Cl2 N4 O5 S

Double bond geometry as shown.

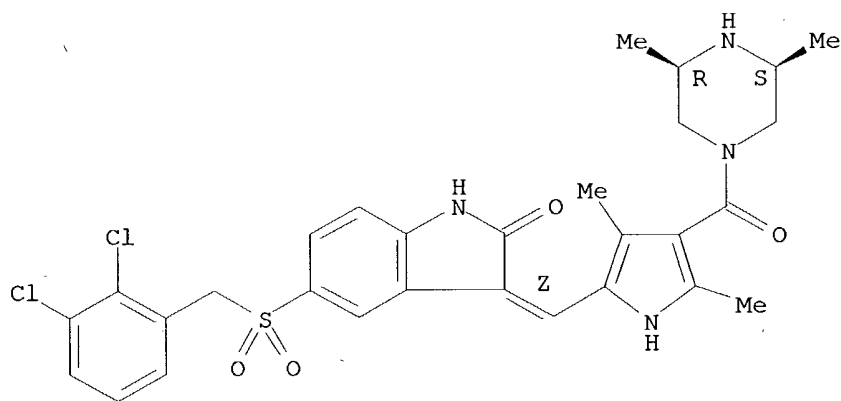


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[(Z)-[5-[[[2,3-dichlorophenyl)methyl]sulfonyl]-1,2-
dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-
yl]carbonyl]-3,5-dimethyl-, (3R,5S)- (9CI)
MF C29 H30 Cl2 N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.

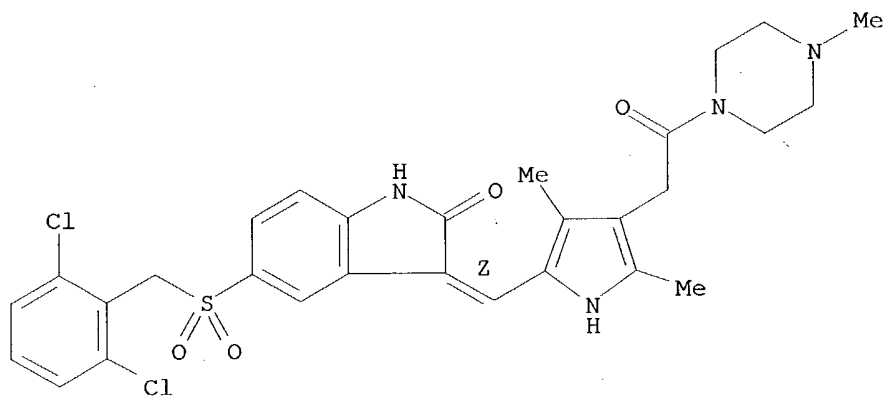


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[(Z)-[5-[[[(2,6-dichlorophenyl)methyl]sulfonyl]-1,2-
dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-
yl]acetyl]-4-methyl- (9CI)
MF C29 H30 Cl2 N4 O4 S

Double bond geometry as shown.

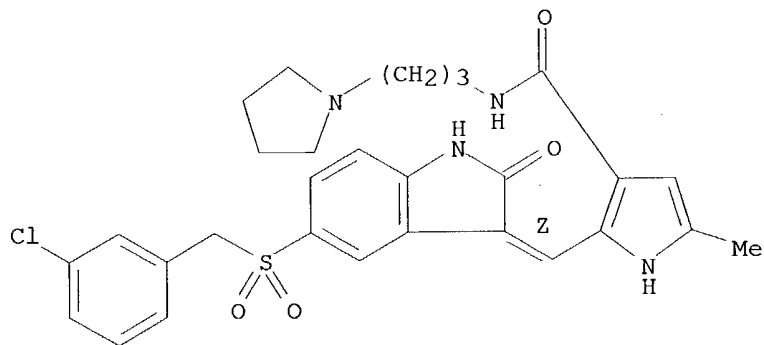


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, 2-[(Z)-[5-[[3-chlorophenyl)methyl]sulfonyl]-1,2-
dihydro-2-oxo-3H-indol-3-ylidene]methyl]-5-methyl-N-[3-(1-
pyrrolidinyl)propyl]- (9CI)
MF C29 H31 Cl N4 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

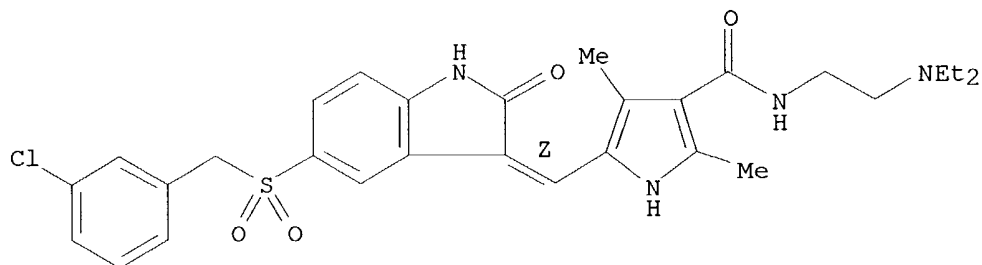
V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Pyrrole-3-carboxamide, 5-[(Z)-[5-[[[(3-chlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-N-[2-(diethylamino)ethyl]-2,4-dimethyl- (9CI)

MF C29 H33 Cl N4 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

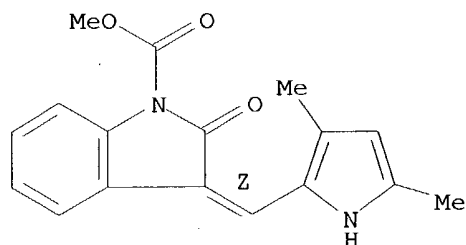
V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Indole-1-carboxylic acid, 3-[(3,5-dimethyl-1H-pyrrol-2-yl)methylene]-
2,3-dihydro-2-oxo-, methyl ester, (3Z)- (9CI)

MF C17 H16 N2 O3

Double bond geometry as shown.



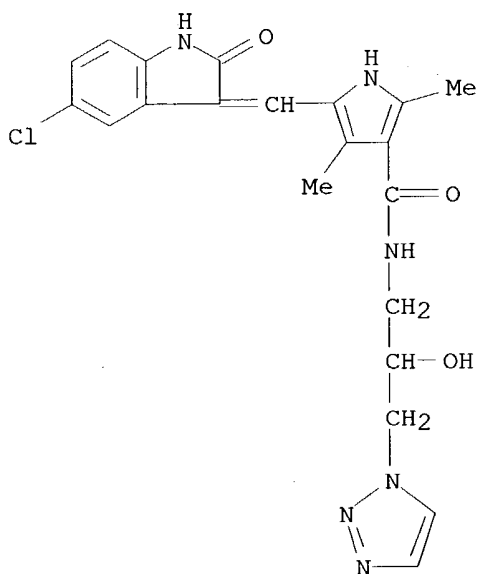
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Pyrrole-3-carboxamide, 5-[(5-chloro-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-N-[2-hydroxy-3-(1H-1,2,3-triazol-1-yl)propyl]-2,4-dimethyl- (9CI)

MF C21 H21 Cl N6 O3



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

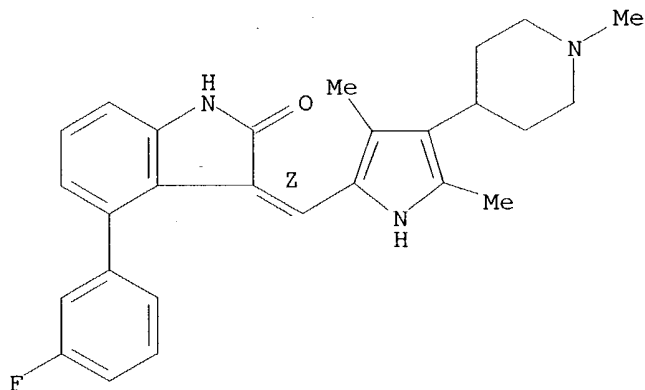
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IN 2H-Indol-2-one, 3-[[3,5-dimethyl-4-(1-methyl-4-piperidinyl)-1H-pyrrol-2-yl]methylene]-4-(3-fluorophenyl)-1,3-dihydro-, (3Z)- (9CI)

MF C27 H28 F N3 O

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

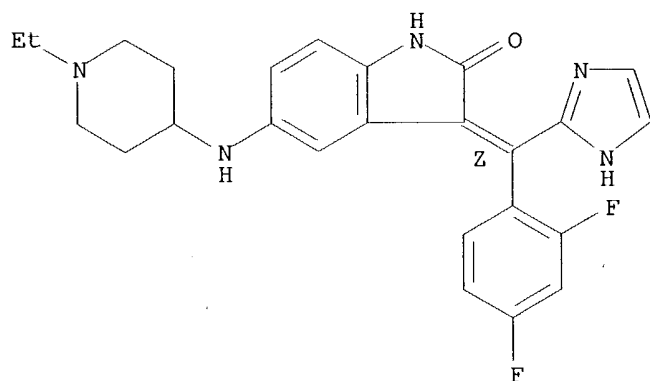
V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 2H-Indol-2-one, 3-[(2,4-difluorophenyl)-1H-imidazol-2-ylmethylene]-5-[(1-ethyl-4-piperidinyl)amino]-1,3-dihydro-, (3Z)- (9CI)

MF C25 H25 F2 N5 O

Double bond geometry as shown.



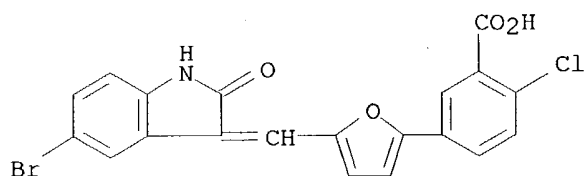
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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IN Benzoic acid, 5-[5-[(5-bromo-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-
2-furanyl]-2-chloro- (9CI)

MF C20 H11 Br Cl N O4



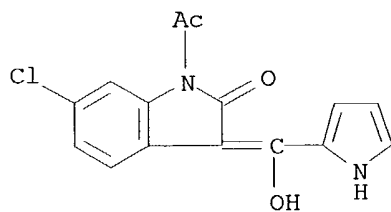
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 2H-Indol-2-one, 1-acetyl-6-chloro-1,3-dihydro-3-(hydroxy-1H-pyrrol-2-ylmethylene)- (9CI)

MF C15 H11 Cl N2 O3

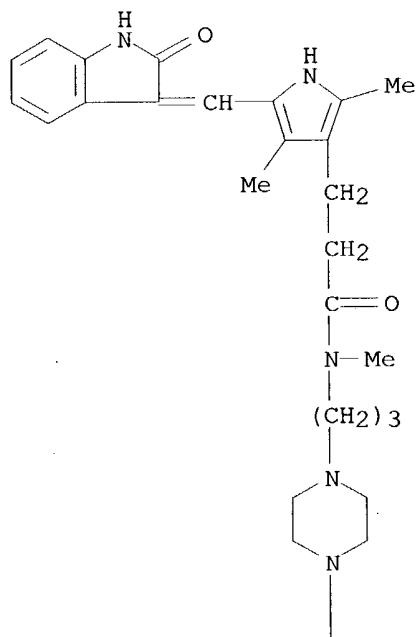


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

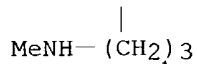
V. Balasubramanian

L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-propanamide, 5-[(1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-
N,2,4-trimethyl-N-[3-[4-[3-(methylamino)propyl]-1-piperazinyl]propyl]-
(9CI)
MF C30 H44 N6 O2
CI COM

PAGE 1-A



PAGE 2-A

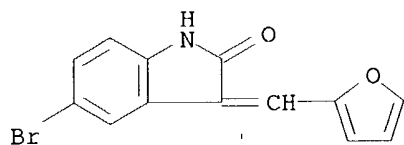


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

10/725,079

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 2H-Indol-2-one, 5-bromo-3-(2-furanylmethylene)-1,3-dihydro- (9CI)
MF C13 H8 Br N O2

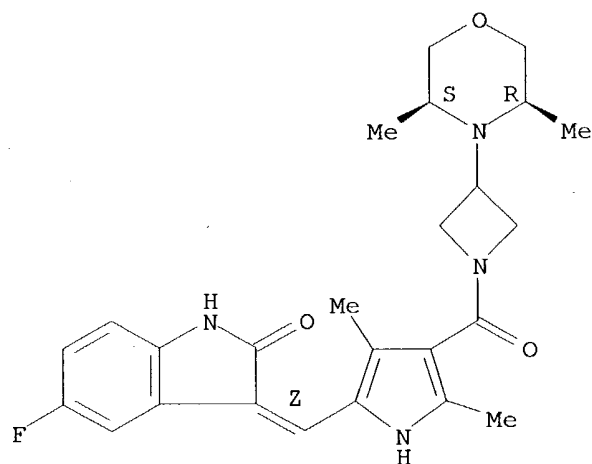


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Azetidine, 3-[(3R,5S)-3,5-dimethyl-4-morpholinyl]-1-[[5-[(Z)-(5-fluoro-1,2-
dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2,4-dimethyl-1H-pyrrol-3-
yl]carbonyl]-, rel- (9CI)
MF C25 H29 F N4 O3

Relative stereochemistry.
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

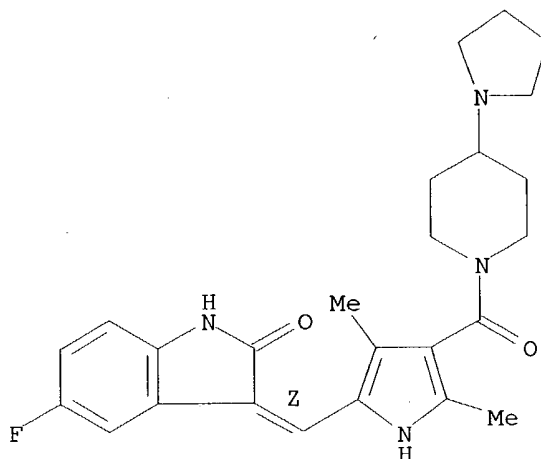
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Piperidine, 1-[[5-[(Z)-(5-fluoro-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-4-(1-pyrrolidinyl)-(9CI)

MF C25 H29 F N4 O2

Double bond geometry as shown.

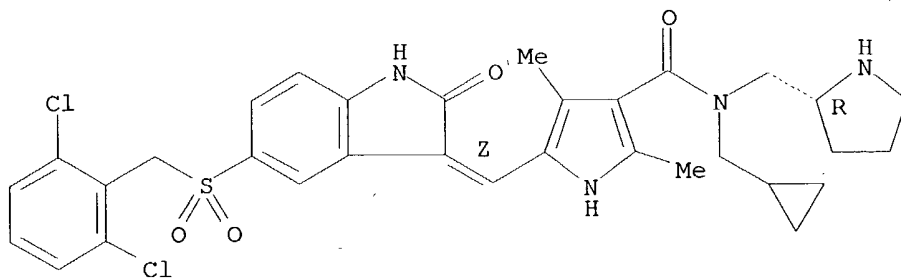


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN 1H-Pyrrole-3-carboxamide, N-(cyclopropylmethyl)-5-[(Z)-[5-[[(2,6-
dichlorophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-
ylidene]methyl]-2,4-dimethyl-N-[(2R)-2-pyrrolidinylmethyl]- (9CI)
MF C32 H34 Cl2 N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

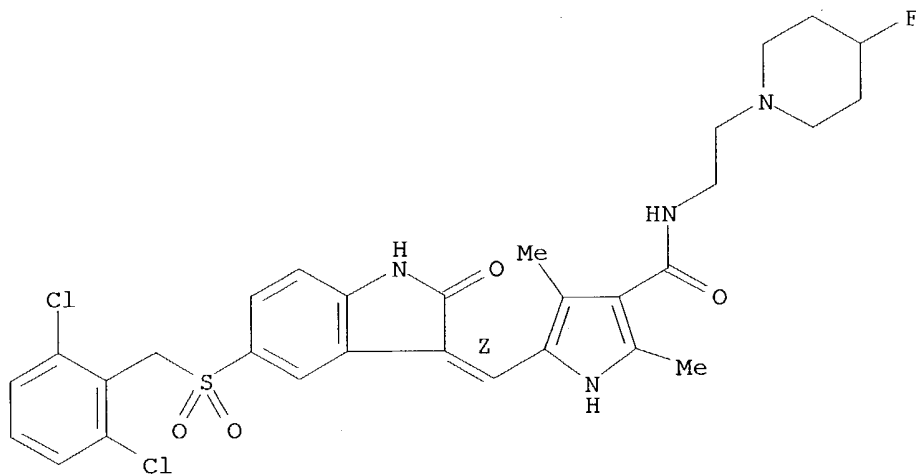
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Pyrrole-3-carboxamide, 5-[(Z)-[5-[(2,6-dichlorophenyl)methyl]sulfonyl]-
1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-N-[2-(4-fluoro-1-
piperidinyl)ethyl]-2,4-dimethyl- (9CI)

MF C30 H31 Cl2 F N4 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

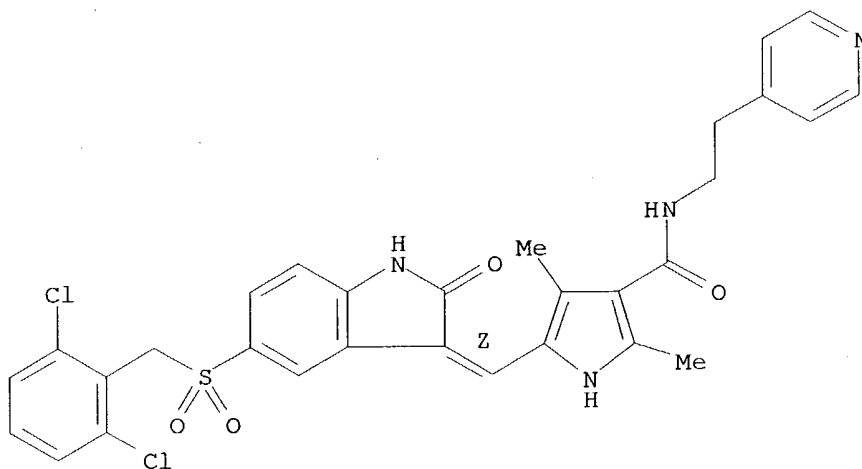
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Pyrrole-3-carboxamide, 5-[(Z)-[5-[[[2,6-dichlorophenyl)methyl]sulfonyl]-
1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-2,4-dimethyl-N-[2-(4-
pyridinyl)ethyl]- (9CI)

MF C30 H26 Cl2 N4 O4 S

Double bond geometry as shown.

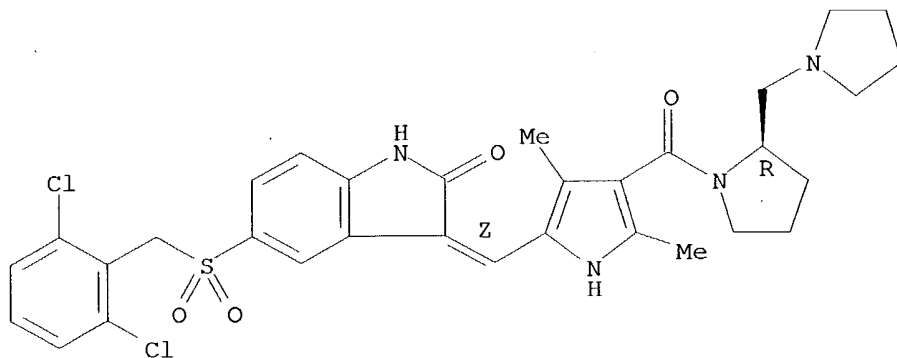


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Pyrrolidine, 1-[[5-[(Z)-[5-[[2,6-dichlorophenyl)methyl]sulfonyl]-1,2-
dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-
yl]carbonyl]-2-(1-pyrrolidinylmethyl)-, (2R)- (9CI)
MF C32 H34 Cl2 N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.

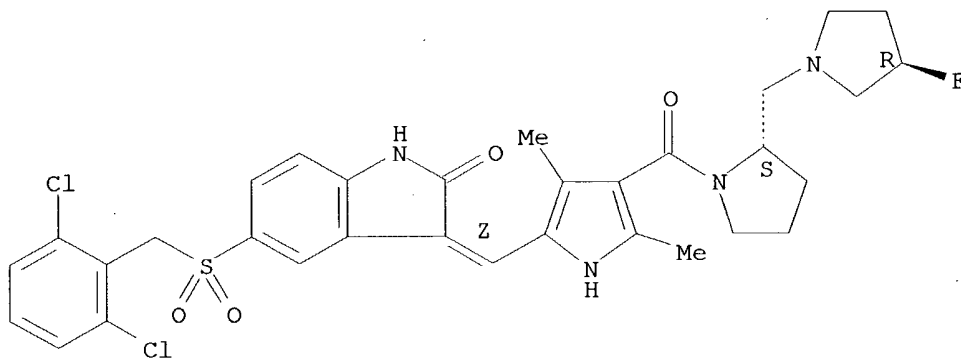


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Pyrrolidine, 1-[[[5-[(Z)-[5-[[[2,6-dichlorophenyl)methyl]sulfonyl]-1,2-
dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-
yl]carbonyl]-2-[[[3R)-3-fluoro-1-pyrrolidinyl]methyl]-, (2S)- (9CI)
MF C32 H33 Cl2 F N4 O4 S

Absolute stereochemistry.
Double bond geometry as shown.

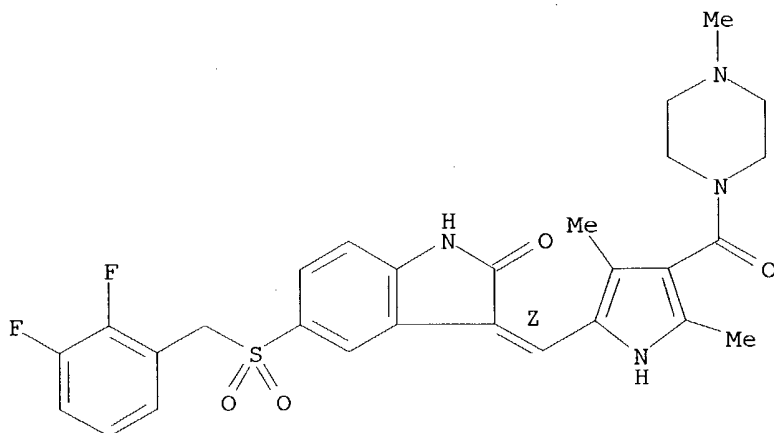


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[(Z)-[5-[[2,3-difluorophenyl)methyl]sulfonyl]-1,2-
dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-
yl]carbonyl]-4-methyl- (9CI)
MF C28 H28 F2 N4 O4 S

Double bond geometry as shown.

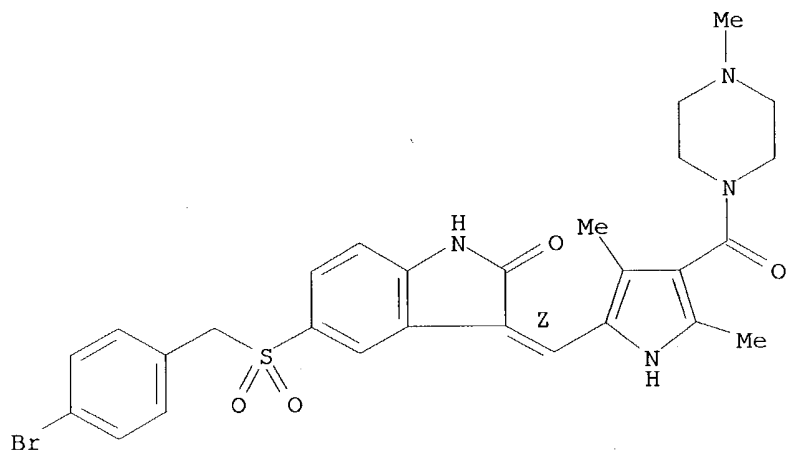


PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Piperazine, 1-[[5-[(Z)-[5-[[4-bromophenyl)methyl]sulfonyl]-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-4-methyl- (9CI)
MF C28 H29 Br N4 O4 S

Double bond geometry as shown.



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

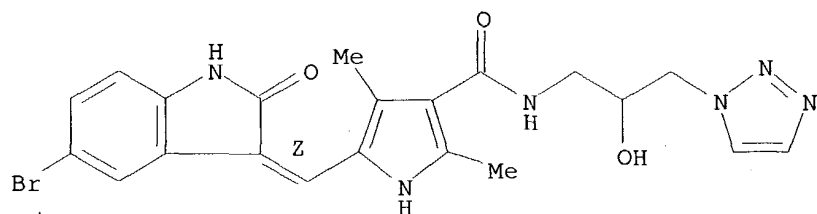
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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Pyrrole-3-carboxamide, 5-[(Z)-(5-bromo-1,2-dihydro-2-oxo-3H-indol-3-ylidene)methyl]-N-[2-hydroxy-3-(1H-1,2,3-triazol-1-yl)propyl]-2,4-dimethyl- (9CI)

MF C21 H21 Br N6 O3

Double bond geometry as shown.



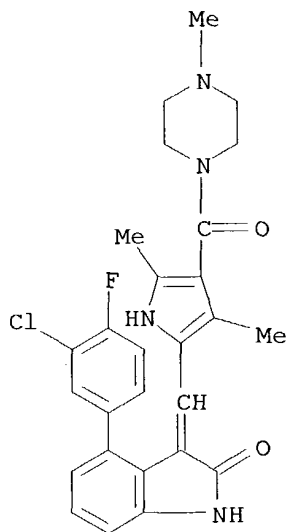
PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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L3 50 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN Piperazine, 1-[[5-[[4-(3-chloro-4-fluorophenyl)-1,2-dihydro-2-oxo-3H-indol-3-ylidene]methyl]-2,4-dimethyl-1H-pyrrol-3-yl]carbonyl]-4-methyl- (9CI)

MF C27 H26 Cl F N4 O2



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

V. Balasubramanian

=> log y

COST IN U.S. DOLLARS

SINCE FILE

ENTRY

TOTAL

SESSION

FULL ESTIMATED COST

0.84

1.05

STN INTERNATIONAL LOGOFF AT 16:39:34 ON 17 AUG 2004